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=> file caplus

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L1 STRUCTURE UPLOADED

=> d

L1 HAS NO ANSWERS

L1 STR

G1 0, S

Structure attributes must be viewed using STN Express query preparation.

=> s 11

REG1stRY INITIATED

Substance data SEARCH and crossover from CAS REGISTRY in progress... Use DISPLAY HITSTR (or FHITSTR) to directly view retrieved structures.

TOh

SAMPLE SEARCH INITIATED 15:48:54 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 63 TO ITERATE

100.0% PROCESSED 63 ITERATIONS 13 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 784 TO 1736
PROJECTED ANSWERS: 44 TO 476

L2 13 SEA SSS SAM L1

L3 10 L2

=> d 1-10 ibib abs hitstr

L3 ANSWER 1 OF 10 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2004:203802 CAPLUS

DOCUMENT NUMBER: 140:235428

TITLE: Preparation of malononitrile compound and use thereof

as pesticides

INVENTOR(S): Okada, Satoshi; Oohira, Daisuke; Otaka, Ken PATENT ASSIGNEE(S): Sumitomo Chemical Company, Limited, Japan

SOURCE: PCT Int. Appl., 104 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT N	K	KIND DATE				APPLICATION NO.					DATE			
WO 20040	WO 2004020399			A1 20040311			WO 2003-JP10726					20030826		
. W:	AE, AG,	AL, A	1, AT,	AU,	AZ,	BA,	BB,	BG,	BR,	BY,	BZ,	CA,	CH,	CN,
	CO, CR,													
•	GM, HR,	HU, I	, IL,	IN,	IS,	KE,	KG,	KR,	KZ,	LC,	LK,	LR,	LS,	LT,
	LU, LV,					•	,				•	•		-
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	KG, KZ,			•							•		-	
	FI, FR,	=		•	•				•		•			-
	BF, BJ,	•	•	•	•	•	•	•	•	•	•	•		•
AU 20032		•				AU 2003-256083				•	20030826			
BR 20030	13964		Ā	2005	0719		BR 2	003-	13964	1		2	0030	826
CN 16785			A	2005	1005		CN 2	003-	8204	24		2	0030	826
JP 20041	43148		4	2004	0520		JP 2	003-	20899	94		2	0030	827
US 20060	04092		1	20060105			US 2005-522764		54		2	0050	201	
PRIORITY APPLN. INFO.:							JP 2	002-	2503	55	7	A 2	0020	B29
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OTUED COIDCE/	el.	M	יייאמם	140.	22541						·	_		

OTHER SOURCE(S): MARPAT 140:235428

GΙ

AΒ The present invention relates to a novel malononitrile compound represented by the formula (I): wherein, R1 represents C1 to C6 alkyl that may be substituted with halogen, C2 to C6 alkenyl that may be substituted with halogen, etc; R2 represents hydrogen atom or C1 to C6 alkyl that may be substituted with halogen; R3 represents hydrogen atom or C1 to C6 alkyl; R4 represents hydrogen atom or C1 to C6 alkyl; R5 represents C1 to C6 alkyl that may be substituted with halogen, C3 to C6 alkenyl that may be substituted with halogen, etc , or R4 and R5 may be combined at their terminal and represent ethylene that may be substituted with C1 to C3 alkyl or trimethylene that may be substituted with C1 to C3 alkyl; and Z1 and Z2, which are the same or different, represent oxygen atom or sulfur atom. Thus, 2-(tert-butoxycarbonylmethyl)-2-allylmalononitrile was prepared by reacting 2-allylmalonitrile with tert-Bu bromoacetate in DMF in the presence of sodium hydride. The malononitrile compound has an efficient pesticidal activity and can control effectively pests such as insect pests, acarine pests, nematode pests and the like.

RN 666738-88-1 CAPLUS

CN Hexanoic acid, 3,3-dicyano-4-methyl-, 1,1-dimethylethyl ester (CA INDEX

RN 666738-93-8 CAPLUS CN Hexanoic acid, 3,3-dicyano-, 1,1-dimethylethyl ester (CA INDEX NAME)

RN 666738-94-9 CAPLUS

CN Heptanoic acid, 3,3-dicyano-2-methyl-, ethyl ester (CA INDEX NAME)

RN 666738-97-2 CAPLUS

CN Heptanoic acid, 3,3-dicyano-2-ethyl-, ethyl ester (CA INDEX NAME)

RN 666739-14-6 CAPLUS

CN Hexanoic acid, 3,3-dicyano-6,6,6-trifluoro-, 3-methoxy-3-methylbutyl ester (CA INDEX NAME)

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L3 ANSWER 2 OF 10 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2000:141061 CAPLUS

DOCUMENT NUMBER: 132:278722

TITLE: Spontaneous addition of active methine compounds to

enol ethers and α , β -unsaturated ketones in

aprotic polar solvent

AUTHOR(S): Yokozawa, Tsutomu; Oishi, Motoi; Tanaka, Yasukazu

CORPORATE SOURCE: Department of Applied Chemistry, Kanagawa University,

Kanagawa-ku Yokohama, 221-8686, Japan

SOURCE: Journal of Organic Chemistry (2000), 65(6), 1895-1897

CODEN: JOCEAH; ISSN: 0022-3263

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal LANGUAGE: English

OTHER SOURCE(S): CASREACT 132:278722

AB Addition of (EtO)2CHCH2CXYCH(CN)2 (I, X = Y = cyano; X = cyano, Y = CO2Me; X = Y = CO2Me) to enol ethers and α, β -unsatd. ketones in DMF at room temp is reported. Thus, reacting I (X = Y = cyano) with H2C:CHOEt gave (EtO)2CHCH2C(CN)2CH(OEt)Me in 63% yield. This reaction illustrates that the electron-withdrawing groups at the β -positions of the active methine group having the ones at the α and β positions were

strongly affected on the acidity of I.

IT 264142-40-7P 264142-41-8P 264142-43-0P

RL: SPN (Synthetic preparation); PREP (Preparation) (addition of methine compds. to enol ethers and α,β -unsatd. ketones)

RN 264142-40-7 CAPLUS

CN Heptanoic acid, 2,3,3-tricyano-2-(2,2-diethoxyethyl)-6-oxo-, methyl ester (CA INDEX NAME)

RN 264142-41-8 CAPLUS

CN Propanedioic acid, (1,1-dicyano-4-oxopentyl)(2,2-diethoxyethyl)-, dimethyl ester (9CI) (CA INDEX NAME)

RN 264142-43-0 CAPLUS

CN Octanoic acid, 2,3,3-tricyano-2-(2,2-diethoxyethyl)-6-oxo-, methyl ester (CA INDEX NAME)

REFERENCE COUNT:

THERE ARE 25 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L3 ANSWER 3 OF 10 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 1999:142376 CAPLUS

DOCUMENT NUMBER: 130:239567

TITLE: Diazaspirononanium salt for use as template for

zeolite synthesis

INVENTOR(S): Kubota, Yoshihiro; Sugi, Yoshihiro

PATENT ASSIGNEE(S): Showa Denko K. K., Japan SOURCE: Jpn. Kokai Tokkyo Koho, 9 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 11060577	A	19990302	JP 1997-220414	19970815
PRIORITY APPLN. INFO.:			JP 1997-220414	19970815

OTHER SOURCE(S): MARPAT 130:239567

AB Claimed template is a salt of substituted 2,7-diazaspiro[4,4]nonanium. Hydrothermal synthesis of a zeolite by bringing a silica source and/or an alumina source into contact with the zeolite is also claimed. ZSM-12 zeolites having crystal size of a major axis $\geq\!50~\mu m$ are also claimed.

IT 77415-69-1P

RL: IMF (Industrial manufacture); RCT (Reactant); PREP (Preparation); RACT (Reactant or reagent)

(reaction of; diazaspirononanium salts as templates for manufacture of ZSM-12 zeolites having large crystal size)

RN 77415-69-1 CAPLUS

CN Pentanedioic acid, 3,3-dicyano-, diethyl ester (9CI) (CA INDEX NAME)

$$\begin{array}{c|cccc} \mathtt{O} & \mathtt{CN} & \mathtt{O} \\ || & | & || \\ \mathtt{EtO-C-CH}_2-\mathtt{C-CH}_2-\mathtt{C-OEt} \\ | & \\ \mathtt{CN} \end{array}$$

L3 ANSWER 4 OF 10 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 1995:900662 CAPLUS

DOCUMENT NUMBER: 124:116317

TITLE: Lanthanum isopropoxide catalyzed addition of activated

nucleophiles to imines

AUTHOR(S): Yamamoto, Yoshinori; Fukui, Hiroyuki; Honda, Yoshihiro

CORPORATE SOURCE: Dept. Chem., Tohoku Univ., Sendai, 980-77, Japan SOURCE: Applied Organometallic Chemistry (1995), 9(5 & 6),

467-71

CODEN: AOCHEX; ISSN: 0268-2605

PUBLISHER: Wiley
DOCUMENT TYPE: Journal
LANGUAGE: English

OTHER SOURCE(S): CASREACT 124:116317

AB The addition of certain activated nucleophiles to activated imines is catalyzed by lanthanum isopropoxide. As activated nucleophiles, methylmaloninitrile and Me 2-cyanopropanoate can be utilized. Imines having an electron-withdrawing group either at the carbon or at the nitrogen atom of the C:N double bond can be used: for example N-toluenesulfonylimines, N-(4-methoxycarbonylphenyl)imines and $\alpha\text{-imino}$ esters.

IT 173006-25-2P

RL: SPN (Synthetic preparation); PREP (Preparation)

(lanthanum isopropoxide catalyzed addition of activated nucleophiles to imines)

RN 173006-25-2 CAPLUS

Butanoic acid, 3,3-dicyano-2-[(1-phenylethyl)amino]-, 5-methyl-2-(1-methyl-CN 1-phenylethyl)cyclohexyl ester, $[1R-[1\alpha[R*(R*)],2\beta,5\alpha]]$ -(CA INDEX NAME)

Absolute stereochemistry.

CAPLUS COPYRIGHT 2008 ACS on STN ANSWER 5 OF 10

ACCESSION NUMBER:

1992:83465 CAPLUS

DOCUMENT NUMBER:

116:83465

TITLE:

The regioselectivity of the ring opening of 1-activated or nonactivated 2-alkoxycarbonyl or 2-cyanoaziridines by carbanions of the dicarbonyl

compounds

AUTHOR (S):

Bouayad, Zoheir; Chanet-Ray, Josette; Ducher, S.;

Vessiere, Roger

CORPORATE SOURCE:

Ec. Natl. Super. Chim. Clermont-Ferrand, Univ. Blaise

Pascal, Aubiere, 63177, Fr.

SOURCE:

Journal of Heterocyclic Chemistry (1991), 28(7),

1757-67

CODEN: JHTCAD; ISSN: 0022-152X

DOCUMENT TYPE:

Journal LANGUAGE: English

GI

Aziridines, e.g. I, reacted with carbanions of dicarbonyl compds., e.g. AB RO2CCH2CO2R (R = Me, Et, CHMe2), to give ring opened products and/or ring enlarged products, e.g. (RO2C) 2CHCH2CH (NHBz) CO2CHMe2, (RO2C)2CHCH(CO2CHMe2)CH2NHBz, and pyrrole II. The regioselectivity depends on several factors. The Ph group on C-3 favors C-3-N bond cleavage, whereas C-2-N bond cleavage is predominant with C-3 substituted or C-2-H aziridines. Cyanoaziridines are predominantly cleaved at C-3-N.

138478-35-0P IT

RL: SPN (Synthetic preparation); PREP (Preparation)

(preparation of)

RN 138478-35-0 CAPLUS

CN Propanoic acid, 2-[(benzoylamino)methyl]-3,3-dicyano-, 1-methylethyl ester (CA INDEX NAME)

L3 ANSWER 6 OF 10 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 1990:497432 CAPLUS

DOCUMENT NUMBER: 113:97432

TITLE: Quinolone antibacterial agents substituted at the

7-position with spiroamines. Synthesis and

structure-activity relationships

AUTHOR(S): Culbertson, Townley P.; Sanchez, Joseph P.; Gambino,

Laura; Sesnie, Josephine A.

CORPORATE SOURCE: Parke-Davis Pharm. Res. Div., Warner-Lambert Co., Ann

Arbor, MI, 48105, USA

SOURCE: Journal of Medicinal Chemistry (1990), 33(8), 2270-5

CODEN: JMCMAR; ISSN: 0022-2623

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 113:97432

GI

AB Fluoroquinolone antibacterials having the 7-position (10-position of pyridobenzoxazines) substituted with 2,7-diazaspiro[4.4]nonane, 1,7-diazaspiro[4.4]nonane, or 2,8-diazaspiro[5.5]undecane (e.g. I (X = CF, CH, N) were prepared and their biol. activities were compared with piperazine and pyrrolidine substituted analogs. Most exhibited potent Gram-pos. and Gram-neg. activity, especially when side chain was N-alkylated. Thus, the quinolinecarboxylic acid II was treated with 2-methyl-2,7-diazaspiro[4.4]nonane to give I (X = CH).

IT 77415-69-1

RL: RCT (Reactant); RACT (Reactant or reagent)
 (reductive cyclization of)

RN 77415-69-1 CAPLUS

CN Pentanedioic acid, 3,3-dicyano-, diethyl ester (9CI) (CA INDEX NAME)

15/01/2008

TOh

ANSWER 7 OF 10 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER:

1986:496793 CAPLUS

DOCUMENT NUMBER:

105:96793

CORPORATE SOURCE:

ORIGINAL REFERENCE NO.: 105:15633a,15636a

TITLE:

Zwitterionic tetramethylenes as the common

intermediates in the cycloaddition and polymerization reactions of N-vinylcarbazole with electrophilic tetrasubstituted ethylenes: a new explanation for

charge-transfer initiation

AUTHOR(S):

Gotoh, Tetsuya; Padias, Anne Buyle; Hall, H. K., Jr. Chem. Dep., Univ. Arizona, Tucson, AZ, 85721, USA

SOURCE:

Journal of the American Chemical Society (1986),

108(16), 4920-31

CODEN: JACSAT; ISSN: 0002-7863

DOCUMENT TYPE: LANGUAGE:

Journal English

OTHER SOURCE(S):

CASREACT 105:96793

The reactions of N-vinylcarbazole (I) with electrophilic tetrasubstituted ethylenes were examples of reactions whose outcomes are manipulated by changes in concentration, structure, and working procedure to form either small mols. (cyclobutanes, 1-butenes) or poly(vinylcarbazole). Equivalent concns. and evaporative workup (organic chemists' conditions) lead to small mols.; a large excess of I and precipitative workup give polymer. The mechanism involves gauche and trans zwitterionic tetramethylenes as intermediates. The former gives cyclobutane reversibly. The latter gives 1-butenes intramol. or adds monomers to form cyclohexanes or eventually polymer. The organic chemical and polymer chemical are unified on this basis. Extensive stereochem. and kinetic support for these propositions is given. Two other proposed mechanisms for these charge-transfer initiations are excluded.

IT 96735-90-9P

> RL: SPN (Synthetic preparation); PREP (Preparation) (preparation of)

RN 96735-90-9 CAPLUS

Propanedioic acid, [2-(9H-carbazol-9-yl)ethenyl](dicyanomethyl)-, dimethyl ester, (E) - (9CI) (CA INDEX NAME)

Double bond geometry as shown.

L3 ANSWER 8 OF 10 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 1985:406747 CAPLUS

DOCUMENT NUMBER: 103:6747

ORIGINAL REFERENCE NO.: 103:1225a,1228a

TITLE: Zwitterionic tetramethylene intermediates: a new

interpretation for "charge-transfer" initiation

AUTHOR(S): Hall, H. K., Jr.; Gotoh, T.

CORPORATE SOURCE: Dep. Chem., Univ. Arizona, Tucson, AZ, 85721, USA

SOURCE: Polymer Preprints (American Chemical Society, Division

of Polymer Chemistry) (1985), 26(1), 34-5

CODEN: ACPPAY; ISSN: 0032-3934

DOCUMENT TYPE: Journal LANGUAGE: English

AB Investigation of the initiation mechanism in polymerization of N-vinylcarbazole (I) [1484-13-5] in the presence of tetracyanoethylene [670-54-2] or di-Me 2,2-dicyanoethylene-1,1-dicarboxylate [82849-49-8] showed that neither the I-cyano compound charge transfer complexes nor the ion-radical pairs formed from them initiated polymerization. The initiating species was the gauche or trans tetramethylene zwitterion formed as an intermediate from the charge-transfer complex. This finding indicated that cyclobutanes initiated vinyl polymerization. The mechanism and the kinetics of the zwitterionic initiation were discussed.

IT 96735-90-9

RL: CAT (Catalyst use); USES (Uses)
 (catalysts, for vinylcarbazole polymerization)

RN 96735-90-9 CAPLUS

CN Propanedioic acid, [2-(9H-carbazol-9-yl)ethenyl](dicyanomethyl)-, dimethyl

ester, (E) - (9CI) (CA INDEX NAME)

Double bond geometry as shown.

L3 ANSWER 9 OF 10 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER:

1981:442180 CAPLUS

DOCUMENT NUMBER:

95:42180

ORIGINAL REFERENCE NO.:

95:7221a,7224a

TITLE:

Absolute configuration of 2,7-diazaspiro[4,4] nonane.

A reassignment

AUTHOR (S):

Overberger, C. G.; Wang, David Wei; Hill, Richard K.;

Krow, Grant R.; Ladner, David W.

CORPORATE SOURCE:

Macromol. Res. Cent., Univ. Michigan, Ann Arbor, MI,

48109, USA

SOURCE:

Journal of Organic Chemistry (1981), 46(13), 2757-64

CODEN: JOCEAH; ISSN: 0022-3263

DOCUMENT TYPE:

Journal

LANGUAGE:

English

OTHER SOURCE(S):

CASREACT 95:42180

GI

The absolute configuration of the axially dissym. spirane 2,7-diazaspiro[4,4] nonane (I), was elucidated as (R)-(-),(S)-(+) in CHCl3 by synthesis of both enantiomers from the centrodissym. intermediate II; the configuration of (R)-(-)-II was correlated with that of (S)-HO2CCMeEtCH2CO2H through the substituted pyrrolidine III. The configuration thus established for the sulfonamide derivative IV is opposite to that derived earlier (Krow, G. and Hill, R. K., 1968). The source of

the original error lies in the preparation of spiroimide V, which is accompanied by almost total racemization when carried out at high temps. A more direct, efficient synthesis of I is described, followed by resolution with dinitrodiphenic acid to give the optically pure enantiomers. Lowe's rule predicts correctly the absolute configurations of several I derivs. but not that of I itself.

IT 77415-69-1P

> RL: SPN (Synthetic preparation); PREP (Preparation) (preparation and diazaspirononane derivative from)

RN 77415-69-1 CAPLUS

Pentanedioic acid, 3,3-dicyano-, diethyl ester (9CI) (CA INDEX NAME) CN

$$\begin{array}{c|cccc} \mathtt{O} & \mathtt{CN} & \mathtt{O} \\ \parallel & \parallel & \parallel \\ \mathtt{EtO-C-CH}_2-\mathtt{C-CH}_2-\mathtt{C-OEt} \\ \parallel & \parallel & \parallel \\ \mathtt{CN} & \end{array}$$

ANSWER 10 OF 10 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 1978:50444 CAPLUS

DOCUMENT NUMBER: 88:50444

ORIGINAL REFERENCE NO.: 88:7949a,7952a

The chemistry of 2-oxopropanedinitrile (carbonyl TITLE:

cyanide); XIX. The ene synthesis using

2-oxopropanedinitrile and 1,3-dicarbonyl compounds

Kociolek, K.; Leplawy, M. T. AUTHOR (S):

Inst. Org. Chem., Tech. Univ. Lodz, Lodz, Pol.
Synthesis (1977), (11), 778-80 CORPORATE SOURCE:

SOURCE: CODEN: SYNTBF; ISSN: 0039-7881

DOCUMENT TYPE: Journal LANGUAGE: English

OTHER SOURCE(S): CASREACT 88:50444

Reaction of CO(CN)2 with RCOCH2COR1 (I; R = R1 = Ph, 2,4,6-Cl3C6H2, Me; R = Me, F3C, R1 = Ph) in ether at 0° was complete in 1 h and gave RCOCH(COR1)C(CN)2OH (II; R and R1 as before) in 100% yield. Reaction of CO(CN)2 with I (R = R1 = OEt) at room temperature required 20 days and gave II in 43-66% yield.

IT 65305-78-4P

> RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation and reaction of, with aniline)

RN65305-78-4 CAPLUS

Propanedioic acid, (dicyanohydroxymethyl)-, diethyl ester (9CI) (CA INDEX CN NAME)

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US Patents Full-Text Database

US OCR Full-Text Database

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<u>L3</u>	malononitril\$9.TI.	331	<u>L3</u>					
<u>L2</u>	L1 AND (514/\$ OR 558/\$)	51	<u>L2</u>					
<u>L1</u>	malononitril\$9 and pesticid\$9	168	<u>L1</u>					

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Search Results - Record(s) 1 through 10 of 10 returned.

☐ 1. Document ID: US 20070117854 A1

L4: Entry 1 of 10

File: PGPB

May 24, 2007 =

PGPUB-DOCUMENT-NUMBER: 20070117854

PGPUB-FILING-TYPE:

DOCUMENT-IDENTIFIER: US 20070117854 A1

TITLE: Malononitrile compounds and use thereof

PUBLICATION-DATE: May 24, 2007

INVENTOR-INFORMATION:

NAME

CITY

STATE

COUNTRY

Mitsudera; Hiromasa

Toyonaka-shi

JP

US-CL-CURRENT: 514/383; 548/267.4

Full Title Citation Front Review Classification Date Reference Sequences Attachments Claims KMC Draw De

☐ 2. Document ID: US 20060004092 A1

L4: Entry 2 of 10

File: PGPB

Jan 5, 2006

PGPUB-DOCUMENT-NUMBER: 20060004092

PGPUB-FILING-TYPE:

DOCUMENT-IDENTIFIER: US 20060004092 A1

TITLE: Malononitrile compound and use thereof pesticides

PUBLICATION-DATE: January 5, 2006

INVENTOR-INFORMATION:

NAME CITY

Takarazuka-shi

STATE

COUNTRY

Okada; Satoshi

Tanarazuna-Siir

JP

Oohra; Daisuke

Toyonaka-shi

JP

Otaka; Ken

Iwaki-shi

JP

US-CL-CURRENT: <u>514/521</u>; <u>558/441</u>

Full Title Citation Front Review Classification Date Reference Sequences Attachments Claims KMC Draw. Dr

7. J. Document ID: US 20050209323 A1

L4: Entry 3 of 10

File: PGPB

Sep 22, 2005

JP

PGPUB-DOCUMENT-NUMBER: 20050209323

PGPUB-FILING-TYPE: new

DOCUMENT-IDENTIFIER: US 20050209323 A1

TITLE: Malononitrile compounds and their use as pesticides

PUBLICATION-DATE: September 22, 2005

INVENTOR-INFORMATION:

COUNTRY NAME CITY STATE

Osaka JΡ Otaka, Ken Oohira, Daisuke Osaka JP Okada, Satoshi Takarazuka-shi

US-CL-CURRENT: 514/520; 558/388

Full Title Citation Front Review Classification Date Reference Sequences Attachments Claims KMC

4. Document ID: US 20050176784 A1

L4: Entry 4 of 10 File: PGPB Aug 11, 2005

PGPUB-DOCUMENT-NUMBER: 20050176784

PGPUB-FILING-TYPE: new

DOCUMENT-IDENTIFIER: US 20050176784 A1

TITLE: Malononitrile compounds and their use as pesticides

PUBLICATION-DATE: August 11, 2005

INVENTOR-INFORMATION:

NAME CITY STATE COUNTRY

Iwaki-shi Otaka, Ken JP J₽ Oohira, Daisuke Toyonaka-shi Takaoka, Daisuke JP Toyonaka-shi

US-CL-CURRENT: <u>514/357</u>; <u>546/320</u>

Full Title Citation Front Review Classification Date Reference Sequences Attachments Claims KMC Draw. De

5. Document ID: US 20040143007 A1

L4: Entry 5 of 10 File: PGPB Jul 22, 2004

PGPUB-DOCUMENT-NUMBER: 20040143007

PGPUB-FILING-TYPE: new

Record List Display Page 3 of 5

DOCUMENT-IDENTIFIER: US 20040143007 A1

TITLE: Malononitrile compounds and their use as pesticides

PUBLICATION-DATE: July 22, 2004

INVENTOR-INFORMATION:

NAME CITY STATE COUNTRY

Otaka, Ken Toyonaka-shi JP
Oohira, Daisuke Minoo-shi JP
Suzuki, Masaya Nishitokyo-shi JP

US-CL-CURRENT: <u>514/520</u>; <u>558/409</u>

Full	Title	Citation	Front	Review	Classification	Date	Reference	Sequences	Attachments	Claims	KWIC	Draw, De

6. Document ID: US 20040138065 A1

L4: Entry 6 of 10 File: PGPB Jul 15, 2004

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